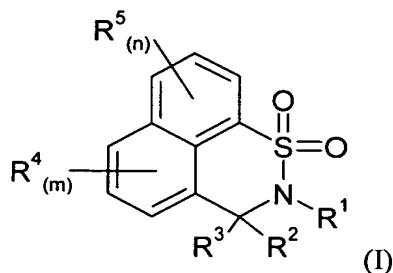


We Claim:

1. A compound of formula (I)



wherein:

R¹ is a group selected from hydrogen, a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, -SO₂H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -CO-C₁-C₆-alkyl, -O, phenyl-C₁-C₄-alkyl, -C₁-C₄-alkyl-NR⁶R⁷, and -C₁-C₄-alkyl-O-C₁-C₄-alkyl, and C₃-C₆-cycloalkyl,

R² and R³, which are identical or different, are each a group selected from hydrogen, a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -NO₂, -SO₂H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -CO-C₁-C₆-alkyl, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -C₁-C₄-alkyl-NR⁶R⁷, and -C₁-C₄-alkyl-O-C₁-C₄-alkyl, and C₃-C₆-cycloalkyl, or

R¹ and R² together are a C₄-C₆-alkylene bridge;

R⁶ and R⁷, which are identical or different, are each hydrogen, C₁-C₄-alkyl, or -CO-C₁-C₄-alkyl;

R⁴, each of which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, phenyl-C₁-C₄-alkyl, halogen, -CN, -NO₂, -SO₂H, -SO₃H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -SO₂-NR⁶R⁷, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-

O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -NR⁶R⁷ and an aryl group optionally mono or polysubstituted by halogen atoms, -NO₂, -SO₂H, or C₁-C₄-alkyl;

R⁵, each of which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, phenyl-C₁-C₄-alkyl, halogen, -CN, -NO₂, -SO₂H, -SO₃H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -SO₂-NR⁶R⁷, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -NR⁶R⁷, and an aryl group optionally mono or polysubstituted by halogen atoms, -NO₂, -SO₂H, or C₁-C₄-alkyl; and

n and m, which are identical or different, are each 0, 1, 2, or 3,

with the proviso that naphtho[1,8-de]-2,3-dihydro-1,1-dioxide-1,2-thiazine is excluded,

or an enantiomer or diastereomer thereof, or a pharmacologically acceptable salt thereof.

2. The compound of formula (I) according to claim 1, wherein:

R¹ is a group selected from hydrogen, a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, -SO₂H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -CO-C₁-C₆-alkyl, -O, -C₁-C₄-alkyl-NR⁷R⁸, and -C₁-C₄-alkyl-O-C₁-C₄-alkyl, benzyl,

R² and R³, which are identical or different, are each a group selected from hydrogen, a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -NO₂, -SO₂H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -CO-C₁-C₆-alkyl, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -C₁-C₄-alkyl-NR⁶R⁷, and -C₁-C₄-alkyl-O-C₁-C₄-alkyl, or

R¹ and R² together are a C₄-C₆-alkylene bridge;

R⁶ and R⁷, which are identical or different, are each hydrogen, C₁-C₄-alkyl, or -CO-C₁-C₂-alkyl, and

R⁴, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -CN, -NO₂, -SO₂H, -SO₃H, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, and -NR⁶R⁷;

R⁵, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -CN, -NO₂, -SO₂H, -SO₃H, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, and -NR⁶R⁷; and

n and m, which are identical or different, are each 0, 1, or 2,

or an enantiomer or diastereomer thereof, or a pharmacologically acceptable salt thereof.

3. The compound of formula (I) according to claim 1, wherein:

R¹ is hydrogen, C₁-C₄-alkyl, or benzyl,

R² and R³, which are identical or different, are each hydrogen or C₁-C₄-alkyl, or

R¹ and R² together are a butylene bridge;

R⁴, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -CN, -NO₂, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, and -NR⁶R⁷;

R⁵, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -CN, -NO₂, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, and -NR⁶R⁷; and

n and m, which are identical or different, are each 0, 1, or 2,

or an enantiomer or diastereomer thereof, or a pharmacologically acceptable salt thereof.

4. The compound of formula (I) according to claim 1, wherein:

R¹, R², R³, which are identical or different, are each hydrogen or C₁-C₄-alkyl;

R⁴, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -NO₂, -O-CO-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -O-C₁-C₆-alkyl, and -NR⁶R⁷;

R⁵, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -NO₂, -O-CO-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -O-C₁-C₆-alkyl, and -NR⁶R⁷; and

n and m, which are identical or different, are each 0, 1, or 2,

or an enantiomer or diastereomer thereof, or a pharmacologically acceptable salt thereof.

5. The compound of formula (I) according to claim 1, wherein:

R¹ is methyl, ethyl, isopropyl, *n*-butyl, or benzyl,

or an enantiomer or diastereomer thereof, or a pharmacologically acceptable salt thereof.

6. The compound of formula (I) according to claim 1, wherein:

R^1 is methyl,

or a pharmacologically acceptable salt thereof.

7. The compound of formula (I) according to claim 1, wherein:

R^1 is methyl;

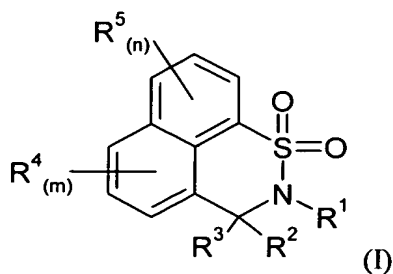
R^2 and R^3 are each hydrogen;

R^4 and R^5 , which are identical or different, are each halogen; and

n and m , which are identical or different, are each 0, 1, or 2,

or a pharmacologically acceptable salt thereof.

8. A compound of general formula (I)



wherein:

R¹ is a group selected from hydrogen, a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, -SO₂H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -CO-C₁-C₆-alkyl, -O, phenyl-C₁-C₄-alkyl, -C₁-C₄-alkyl-NR⁶R⁷, and -C₁-C₄-alkyl-O-C₁-C₄-alkyl, and C₃-C₆-cycloalkyl,

R² and R³, which are identical or different, are each a group selected from hydrogen, a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, halogen, -NO₂, -SO₂H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -CO-C₁-C₆-alkyl, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -C₁-C₄-alkyl-NR⁶R⁷, -C₁-C₄-alkyl-O-, C₁-C₄-alkyl, and C₃-C₆-cycloalkyl, or

R¹ and R² together are a C₄-C₆-alkylene bridge;

R⁶ and R⁷, which are identical or different, are each hydrogen, C₁-C₄-alkyl, or -CO-C₁-C₄-alkyl;

R⁴, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, phenyl-C₁-C₄-alkyl, halogen, -CN, -NO₂, -SO₂H, -SO₃H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -SO₂-NR⁶R⁷, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -NR⁶R⁷, and an aryl group optionally mono or polysubstituted by halogen atoms, -NO₂, -SO₂H, or C₁-C₄-alkyl;

R⁵, which are identical or different, are each a group selected from a C₁-C₆-alkyl group optionally substituted by one or more halogen atoms, phenyl-C₁-C₄-alkyl, halogen, -CN, -NO₂, -SO₂H, -SO₃H, -SO₂-C₁-C₆-alkyl, -SO-C₁-C₆-alkyl, -SO₂-NR⁶R⁷, -COOH, -CO-C₁-C₆-alkyl, -O-CO-C₁-C₄-alkyl, -CO-O-C₁-C₄-alkyl, -O-CO-O-C₁-C₄-alkyl, -CO-NR⁶R⁷, -OH, -O-C₁-C₆-alkyl, -S-C₁-C₆-alkyl, -NR⁶R⁷, and an aryl group optionally mono or polysubstituted by halogen atoms, -NO₂, -SO₂H, or C₁-C₄-alkyl; and

n and m, which are identical or different, are each 0, 1, 2, or 3,

or an enantiomer or diastereomer thereof, or a pharmacologically acceptable salt thereof.